06/04/2008

=>
Uploading 09856396.str

L1 STRUCTURE UPLOADED

=> s l1 sss
SAMPLE SEARCH INITIATED 08:56:49 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 124 TO ITERATE

100.0% PROCESSED 124 ITERATIONS 4 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 1812 TO 3148
PROJECTED ANSWERS: 4 TO 200

L2 4 SEA SSS SAM L1

=> s l1 full
FULL SEARCH INITIATED 08:56:54 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 2532 TO ITERATE

100.0% PROCESSED 2532 ITERATIONS 57 ANSWERS
SEARCH TIME: 00.00.02

L3 57 SEA SSS FUL L1

=> d 1-57 hitstr
'HITSTR' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual
fields or predefined formats. The predefined substance formats
are: (RN = CAS Registry Number)

REG - RN
SAM - Index Name, MF, and structure - no RN
FIDE - All substance data, except sequence data
IDE - FIDE, but only 50 names
SQIDE - IDE, plus sequence data
SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used
SQD - Protein sequence data, includes RN
SQD3 - Same as SQD, but 3-letter amino acid codes are used
SQN - Protein sequence name information, includes RN

CALC - Table of calculated properties
EPROP - Table of experimental properties
PROP - EPROP and CALC

Any CA File format may be combined with any substance format to
obtain CA references citing the substance. The substance formats
must be cited first. The CA File predefined formats are:

ABS -- Abstract
APPS -- Application and Priority Information
BIB -- CA Accession Number, plus Bibliographic Data
CAN -- CA Accession Number
CBIB -- CA Accession Number, plus Bibliographic Data (compressed)
IND -- Index Data
IPC -- International Patent Classification

PATS -- PI, SO
STD -- BIB, IPC, and NCL

IABS --ABS, indented, with text labels
IBIB -- BIB, indented, with text labels
ISTD -- STD format, indented

OBIB -------- AN, plus Bibliographic Data (original)
OIBIB ------- OBIB, indented with text labels

SBIB -------- BIB, no citations
SIBIB ------- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when
it is available.
The MAX format is the same as ALL.
The IALL format is the same as ALL with BIB ABS and IND indented,
with text labels.

For additional information, please consult the following help
messages:

HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
ENTER DISPLAY FORMAT (IDE):str\
'STR\' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

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are: (RN = CAS Registry Number)

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SQD   - Protein sequence data, includes RN
SQD3   - Same as SQD, but 3-letter amino acid codes are used
SQN   - Protein sequence name information, includes RN

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EPROP - Table of experimental properties
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IND -- Index Data
IPC -- International Patent Classification
PATS -- PI, SO
STD -- BIB, IPC, and NCL

IABS --ABS, indented, with text labels

IBIB -- BIB, indented, with text labels
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OBIB -------- AN, plus Bibliographic Data (original)
OIBIB ------- OBIB, indented with text labels

SBIB -------- BIB, no citations
SIBIB ------- IBIB, no citations

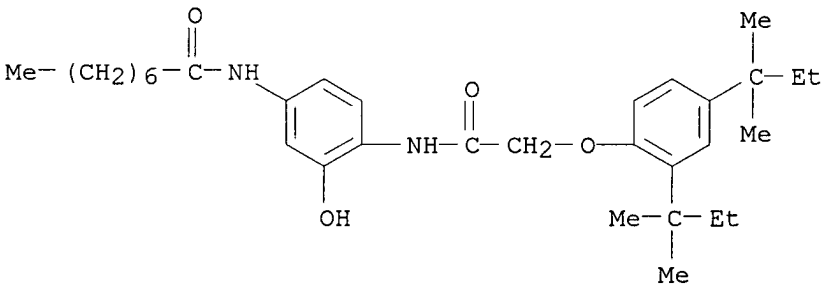
The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.
The MAX format is the same as ALL.
The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

For additional information, please consult the following help messages:

HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
ENTER DISPLAY FORMAT (IDE):end

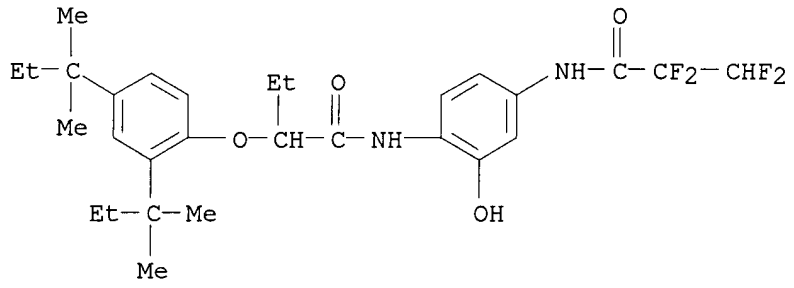
=> d 1-57 sam

L3 ANSWER 1 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Octanamide, N-[4-[[[2,4-bis(1,1-dimethylpropyl)phenoxy]acetyl]amino]-3-hydroxyphenyl]- (9CI)
MF C32 H48 N2 O4



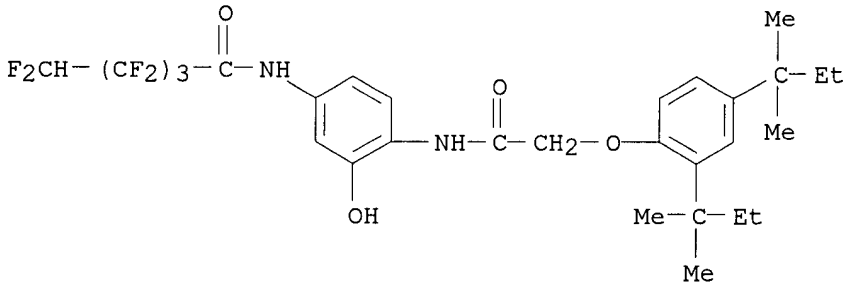
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 2 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Butanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[2-hydroxy-4-[(2,2,3,3-tetrafluoro-1-oxopropyl)amino]phenyl]- (9CI)
MF C29 H38 F4 N2 O4



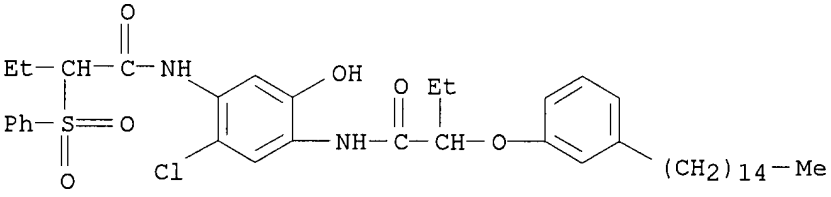
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 3 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Pentanamide, N-[4-[[[2,4-bis(1,1-dimethylpropyl)phenoxy]acetyl]amino]-3-hydroxyphenyl]-2,2,3,3,4,4,5,5-octafluoro- (9CI)
MF C29 H34 F8 N2 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 4 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Butanamide, N-[2-chloro-5-hydroxy-4-[[1-oxo-2-(3-pentadecylphenoxy)butyl]amino]phenyl]-2-(phenylsulfonyl)- (9CI)
MF C41 H57 Cl N2 O6 S



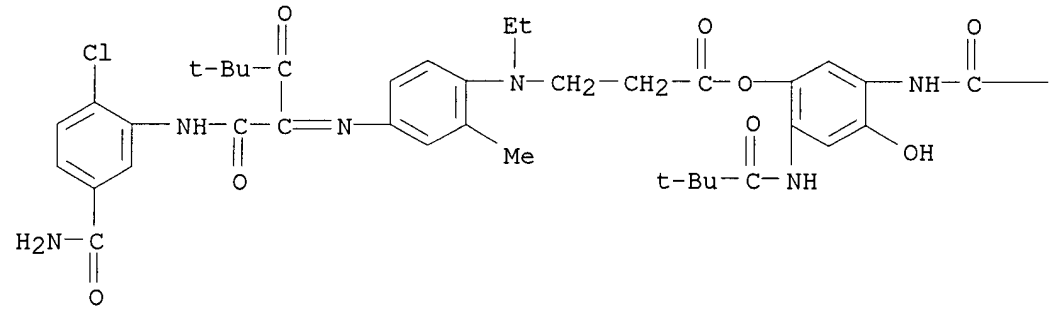
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 5 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN .beta.-Alanine, N-[4-[[1-[[[5-(aminocarbonyl)-2-chlorophenyl]amino]carbonyl]-3,3-dimethyl-2-oxobutylidene]amino]-2-

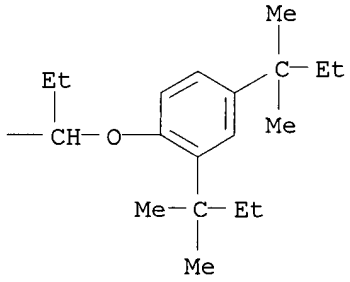
methylphenyl]-N-ethyl-, 5-[[2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-1-oxobutyl]amino]-2-[(2,2-dimethyl-1-oxopropyl)amino]-4-hydroxyphenyl ester (9CI)

MF C57 H75 Cl N6 O9

PAGE 1-A



PAGE 1-B

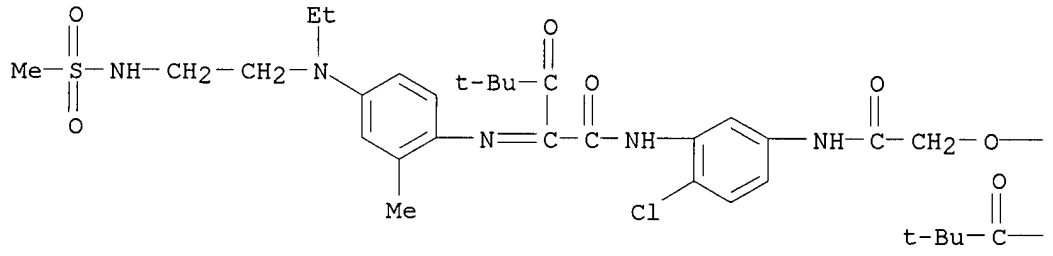


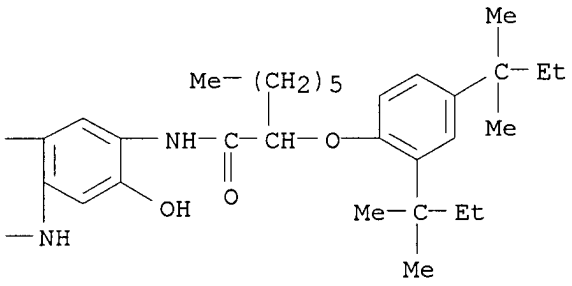
L3 ANSWER 6 OF 57 REGISTRY COPYRIGHT 2002 ACS

IN Octanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[5-[2-[[4-chloro-3-[[2-[[4-[ethyl[2-[(methylsulfonyl)amino]ethyl]amino]-2-methylphenyl]imino]-4,4-dimethyl-1,3-dioxopentyl]amino]phenyl]amino]-2-oxoethoxy]-4-[(2,2-dimethyl-1-oxopropyl)amino]-2-hydroxyphenyl]- (9CI)

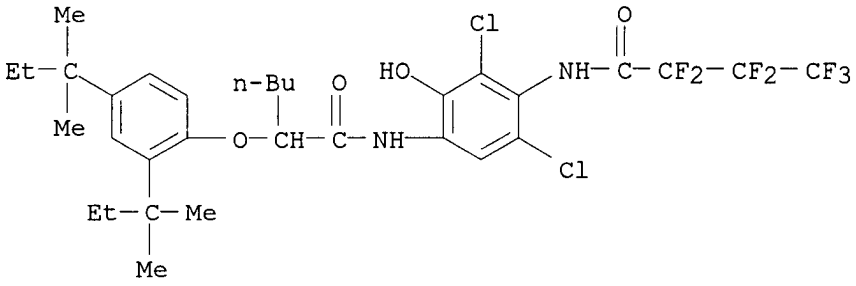
MF C62 H88 Cl N7 O10 S

PAGE 1-A



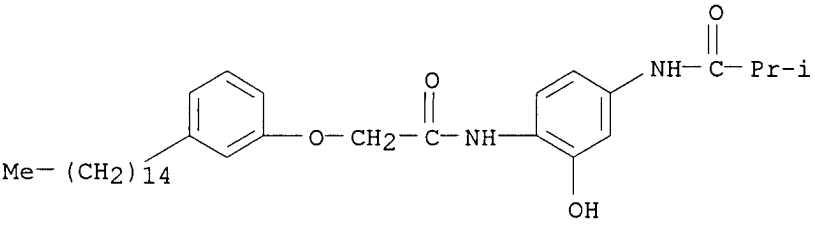


L3 ANSWER 7 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Hexanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[3,5-dichloro-4-[(2,2,3,3,4,4,4-heptafluoro-1-oxobutyl)amino]-2-hydroxyphenyl]- (9CI)
MF C32 H39 Cl2 F7 N2 O4



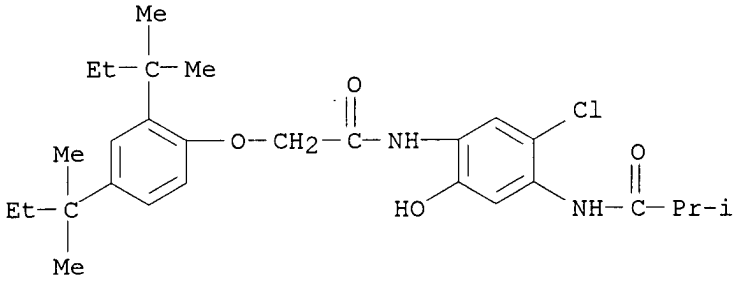
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 8 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Propanamide, N-[3-hydroxy-4-[[(3-pentadecylphenoxy)acetyl]amino]phenyl]-2-methyl- (9CI)
MF C33 H50 N2 O4



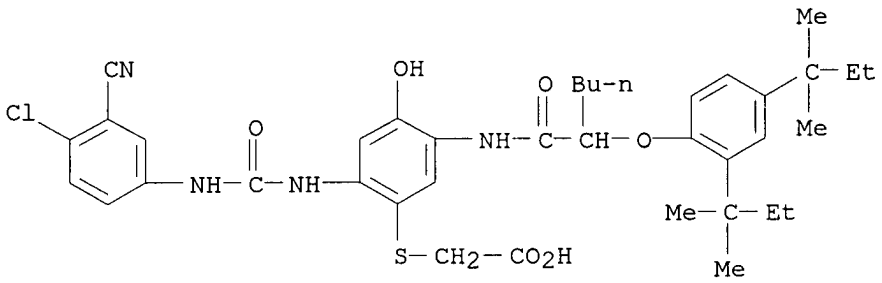
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 9 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Propanamide, N-[4-[[[2,4-bis(1,1-dimethylpropyl)phenoxy]acetyl]amino]-2-chloro-5-hydroxyphenyl]-2-methyl- (9CI)
MF C28 H39 Cl N2 O4



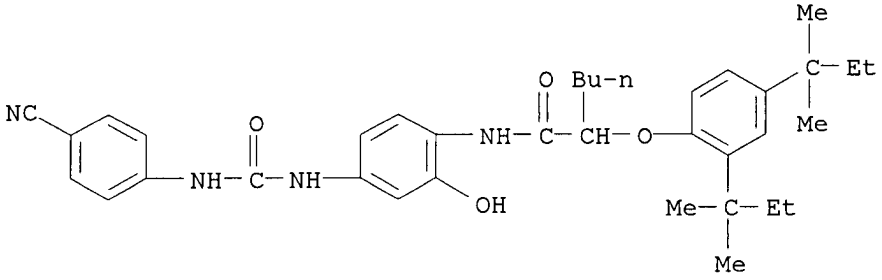
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 10 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Acetic acid, [[5-[[2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-1-oxohexyl]amino]-2-[[[(4-chloro-3-cyanophenyl)amino]carbonyl]amino]-4-hydroxyphenyl]thio]- (9CI)
MF C38 H47 Cl N4 O6 S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

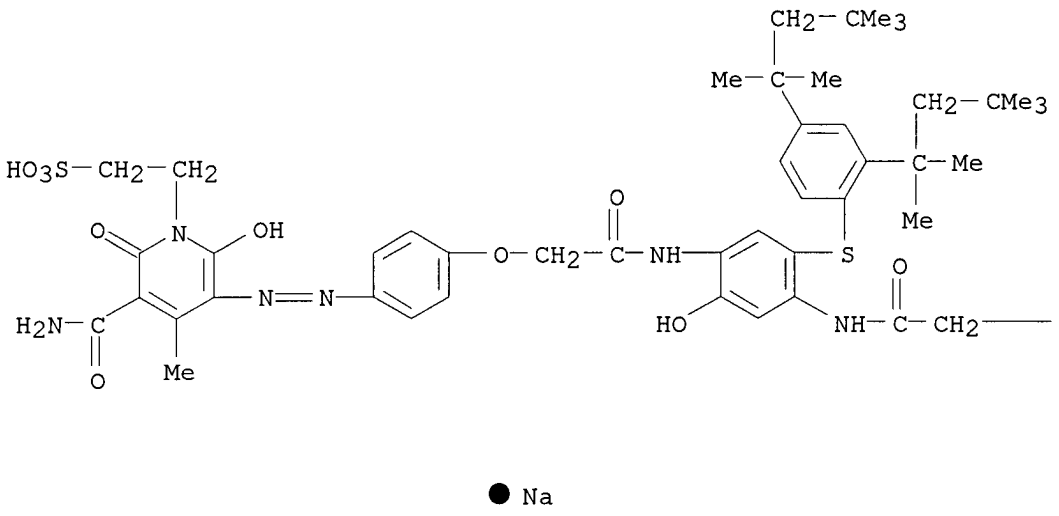
L3 ANSWER 11 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Hexanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[4-[[[(4-cyanophenyl)amino]carbonyl]amino]-2-hydroxyphenyl]- (9CI)
MF C36 H46 N4 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 12 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Butanoic acid, 4-[[4-[[[4-[[5-(aminocarbonyl)-1,6-dihydro-2-hydroxy-4-methyl-6-oxo-1-(2-sulfoethyl)-3-pyridinyl]azo]phenoxy]acetyl]amino]-2-[[2,4-bis(1,1,3,3-tetramethylbutyl)phenyl]thio]-5-hydroxyphenyl]amino]-4-oxo-, monosodium salt (9CI)
MF C49 H64 N6 O12 S2 . Na

PAGE 1-A

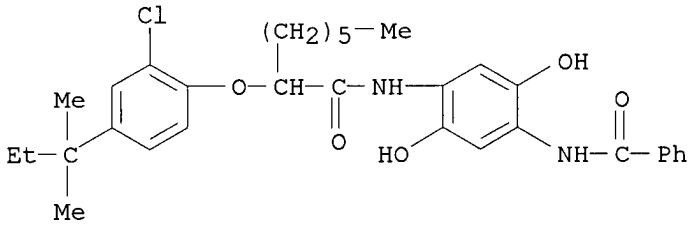


● Na

PAGE 1-B

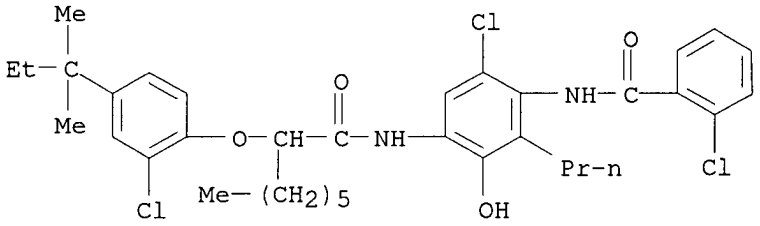
— CH_2 — CO_2H

L3 ANSWER 13 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Benzamide, N-[4-[[2-[2-chloro-4-(1,1-dimethylpropyl)phenoxy]-1-oxooctyl]amino]-2,5-dihydroxyphenyl]- (9CI)
MF C32 H39 Cl N2 O5



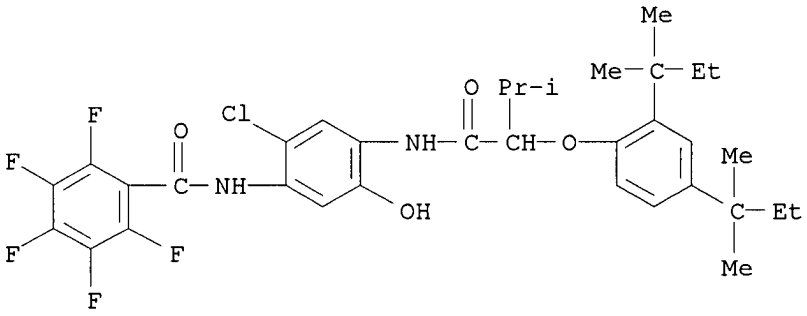
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 14 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Benzamide, 2-chloro-N-[6-chloro-4-[[2-[2-chloro-4-(1,1-dimethylpropyl)phenoxy]-1-oxooctyl]amino]-3-hydroxy-2-propylphenyl]- (9CI)
MF C35 H43 Cl3 N2 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

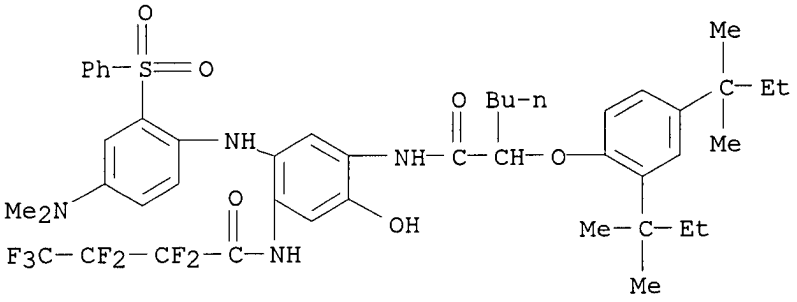
L3 ANSWER 15 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Benzamide, N-[4-[[2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-3-methyl-1-oxobutyl]amino]-2-chloro-5-hydroxyphenyl]-2,3,4,5,6-pentafluoro- (9CI)
MF C34 H38 Cl F5 N2 O4



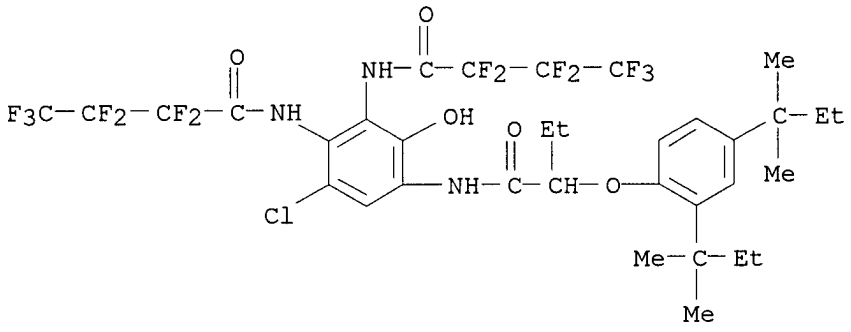
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 16 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Hexanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[5-[[4-(dimethylamino)-2-(phenylsulfonyl)phenyl]amino]-4-[(2,2,3,3,4,4,4-

heptafluoro-1-oxobutyl)amino]-2-hydroxyphenyl]- (9CI)
MF C46 H55 F7 N4 O6 S

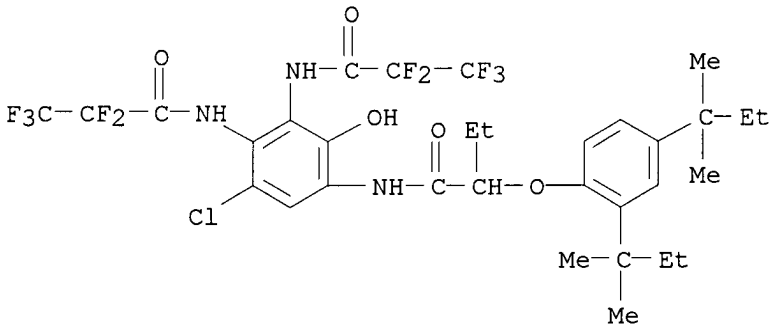


L3 ANSWER 17 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Butanamide, N,N'-[4-[[2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-1-oxobutyl]amino]-6-chloro-3-hydroxy-1,2-phenylene]bis[2,2,3,3,4,4,4-heptafluoro- (9CI)
MF C34 H36 Cl F14 N3 O5



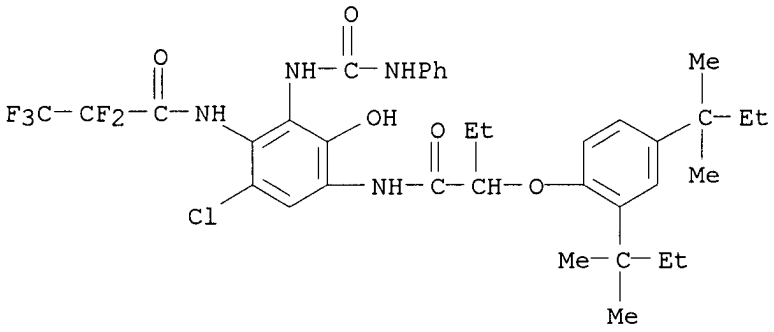
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 18 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Butanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[5-chloro-2-hydroxy-3,4-bis[(2,2,3,3,3-pentafluoro-1-oxopropyl)amino]phenyl]- (9CI)
MF C32 H36 Cl F10 N3 O5



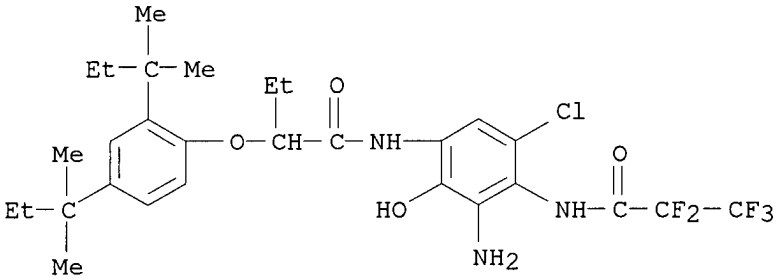
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 19 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Butanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[5-chloro-2-hydroxy-4-[(2,2,3,3,3-pentafluoro-1-oxopropyl)amino]-3-[[(phenylamino)carbonyl]amino]phenyl]- (9CI)
MF C36 H42 Cl F5 N4 O5



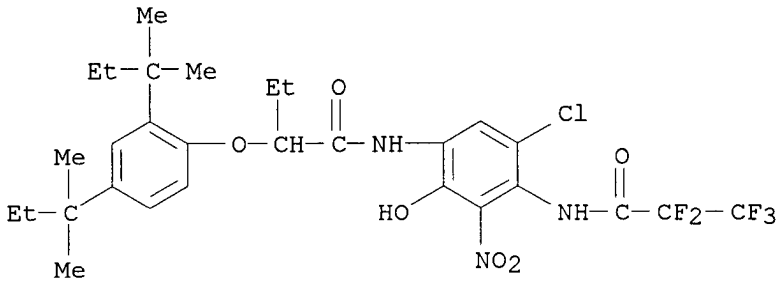
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 20 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Butanamide, N-[3-amino-5-chloro-2-hydroxy-4-[(2,2,3,3,3-pentafluoro-1-oxopropyl)amino]phenyl]-2-[2,4-bis(1,1-dimethylpropyl)phenoxy]- (9CI)
MF C29 H37 Cl F5 N3 O4



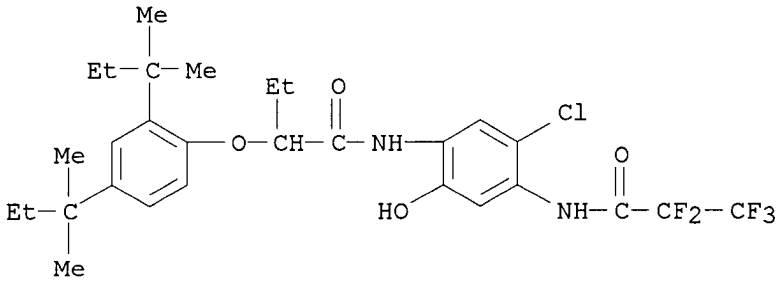
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 21 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Butanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[5-chloro-2-hydroxy-3-nitro-4-[(2,2,3,3,3-pentafluoro-1-oxopropyl)amino]phenyl]- (9CI)
MF C29 H35 Cl F5 N3 O6



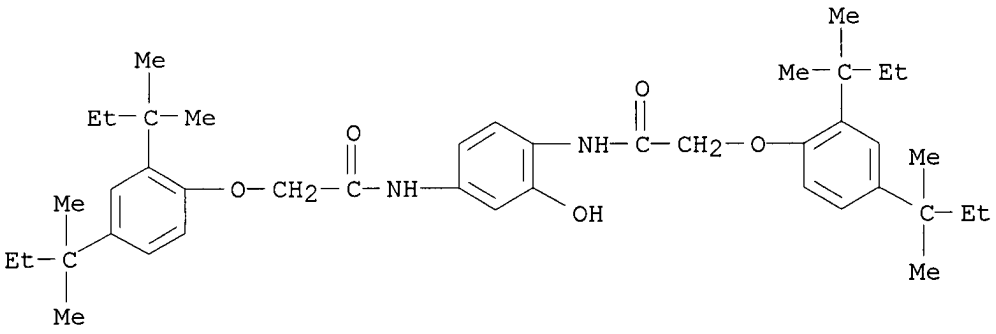
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 22 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Butanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[5-chloro-2-hydroxy-4-[(2,2,3,3,3-pentafluoro-1-oxopropyl)amino]phenyl]- (9CI)
MF C29 H36 Cl F5 N2 O4



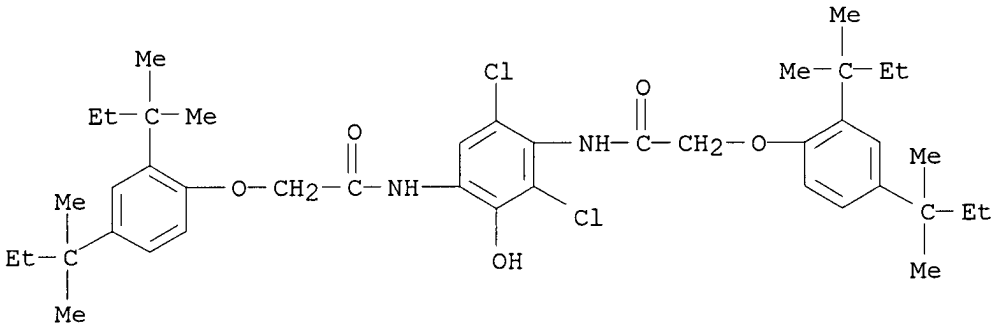
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 23 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Acetamide, N,N'-(hydroxy-p-phenylene)bis[2-(2,4-di-tert-pentylphenoxy)- (6CI)
MF C42 H60 N2 O5



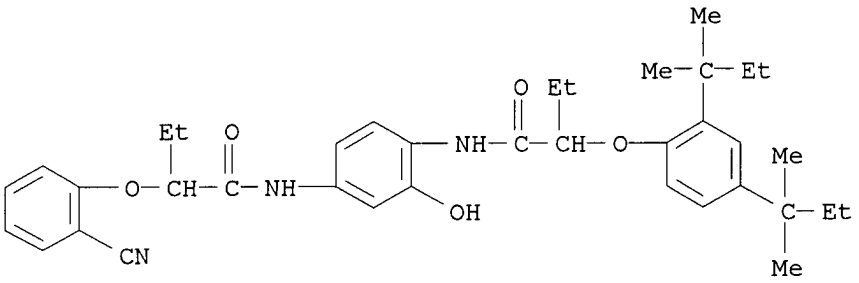
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

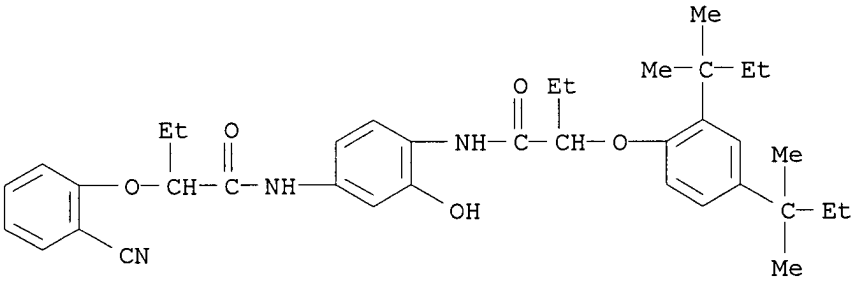
L3 ANSWER 24 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Acetamide, N,N'-(2,6-dichloro-3-hydroxy-p-phenylene)bis[2-(chloro-2,4-di-tert-pentylphenoxy)- (6CI)
MF C42 H56 Cl4 N2 O5
CI IDS



2 (D1—Cl)

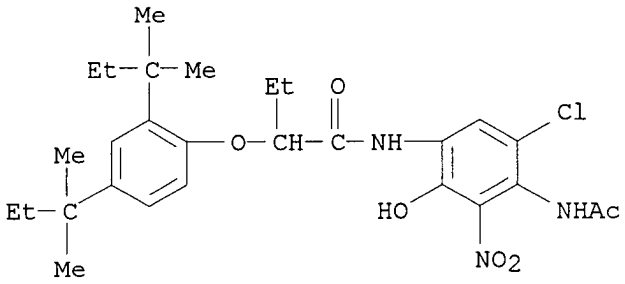
L3 ANSWER 25 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Butanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[4-[[2-(2-cyanophenoxy)-1-oxobutyl]amino]-2-hydroxyphenyl]- (9CI)
MF C37 H47 N3 O5





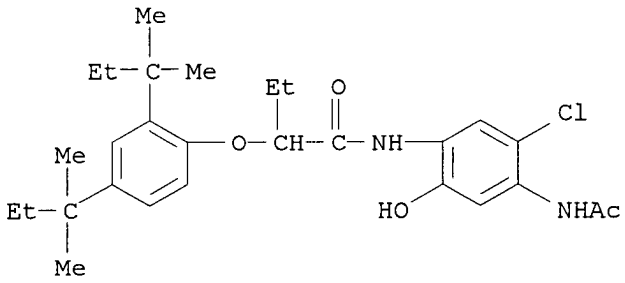
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 26 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Butanamide, N-[4-(acetylamino)-5-chloro-2-hydroxy-3-nitrophenyl]-2-[2,4-bis(1,1-dimethylpropyl)phenoxy]- (9CI)
MF C28 H38 Cl N3 O6



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 27 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Butanamide, N-[4-(acetylamino)-5-chloro-2-hydroxyphenyl]-2-[2,4-bis(1,1-dimethylpropyl)phenoxy]- (9CI)
MF C28 H39 Cl N2 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 28 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Tetradecanamide, N-[3-hydroxy-4-[[[4-[[[3-[[(4-methoxyphenyl)amino]sulfonyl]propyl]sulfonyl]amino]phenoxy]acetyl]amino]phenyl]- (9CI)
MF C38 H54 N4 O9 S2

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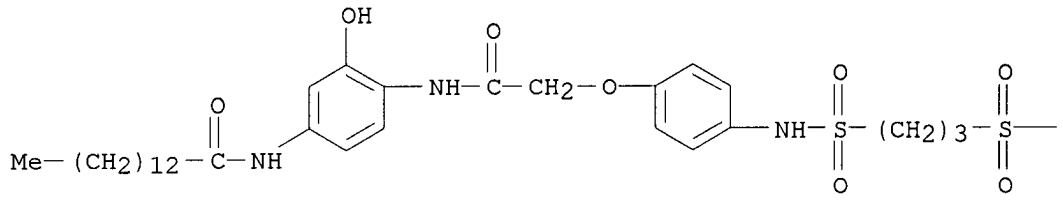

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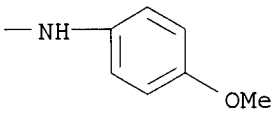
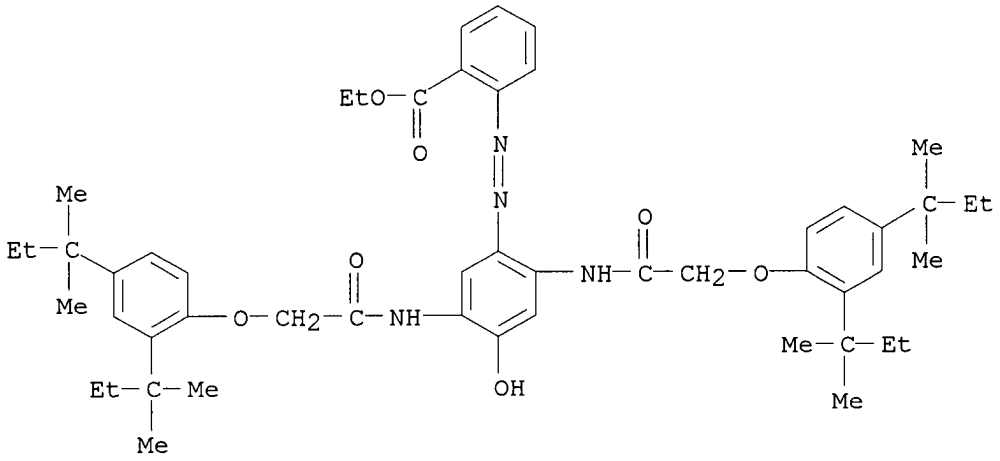


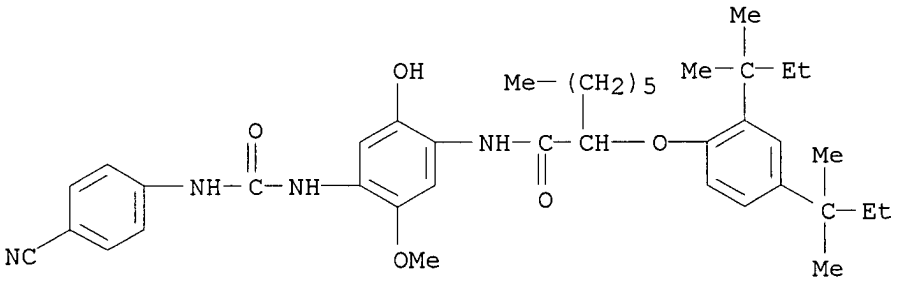
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 29 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Benzoic acid, o-[2,5-bis[2-(2,4-di-tert-pentylphenoxy)acetamido]-4-hydroxyphenylazo]-, ethyl ester (6CI)
MF C51 H68 N4 O7



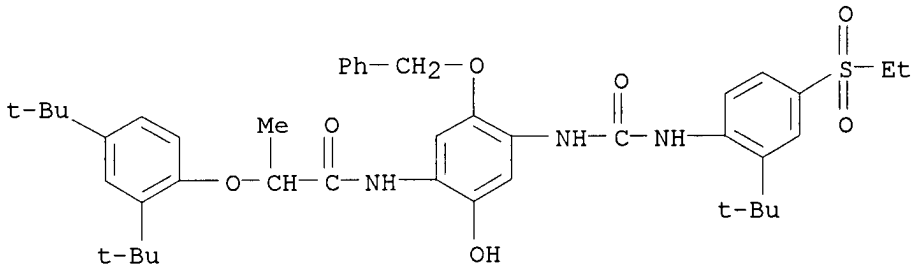
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 30 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Octanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[4-[[[(4-cyanophenyl)amino]carbonyl]amino]-2-hydroxy-5-methoxyphenyl]- (9CI)
MF C39 H52 N4 O5



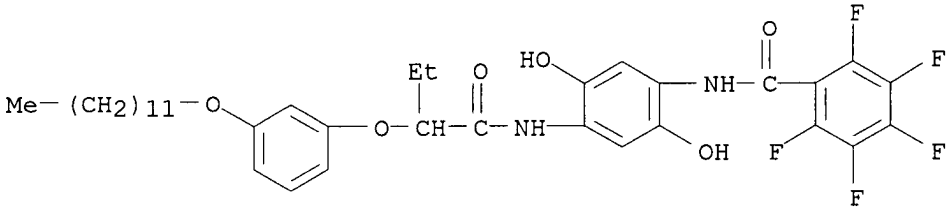
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 31 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Propanamide, 2-[2,4-bis(1,1-dimethylethyl)phenoxy]-N-[4-[[[[2-(1,1-dimethylethyl)-4-(ethylsulfonyl)phenyl]amino]carbonyl]amino]-2-hydroxy-5-(phenylmethoxy)phenyl]- (9CI)
MF C43 H55 N3 O7 S



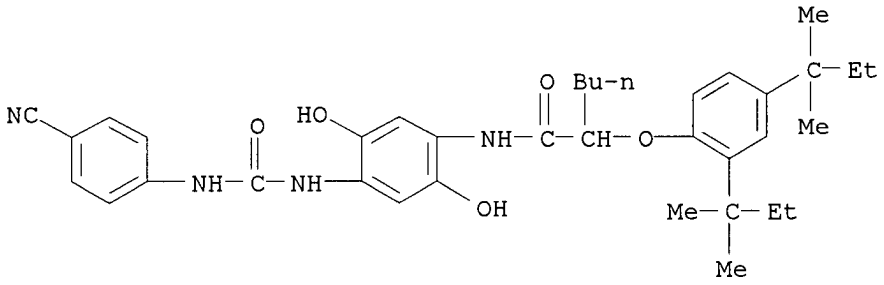
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 32 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Benzamide, N-[4-[[2-[3-(dodecyloxy)phenoxy]-1-oxobutyl]amino]-2,5-dihydroxyphenyl]-2,3,4,5,6-pentafluoro- (9CI)
MF C35 H41 F5 N2 O6



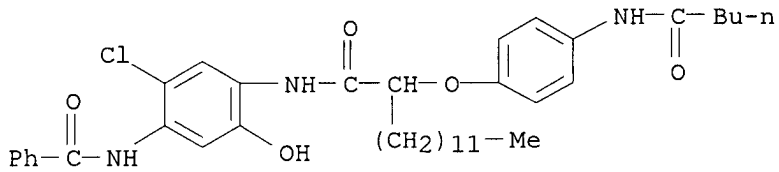
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 33 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Hexanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[4-[[[(4-cyanophenyl)amino]carbonyl]amino]-2,5-dihydroxyphenyl]- (9CI)
MF C36 H46 N4 O5



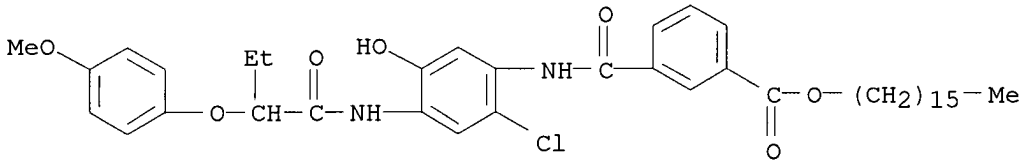
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 34 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Benzamide, N-[2-chloro-5-hydroxy-4-[[1-oxo-2-[4-[(1-oxopentyl)amino]phenoxy]tetradecyl]amino]phenyl]- (9CI)
MF C38 H50 Cl N3 O5



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 35 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Benzoic acid, 3-[[[2-chloro-5-hydroxy-4-[[2-(4-methoxyphenoxy)-1-oxobutyl]amino]phenyl]amino]carbonyl]-, hexadecyl ester (9CI)
MF C41 H55 Cl N2 O7



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 36 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Benzoic acid, 4-[[[5-hydroxy-2-(2-methoxy-2-oxoethoxy)-4-[[2-(3-methoxyphenoxy)-2-methyl-1-oxopropyl]amino]phenyl]amino]carbonyl]-, hexadecyl ester (9CI)
MF C44 H60 N2 O10

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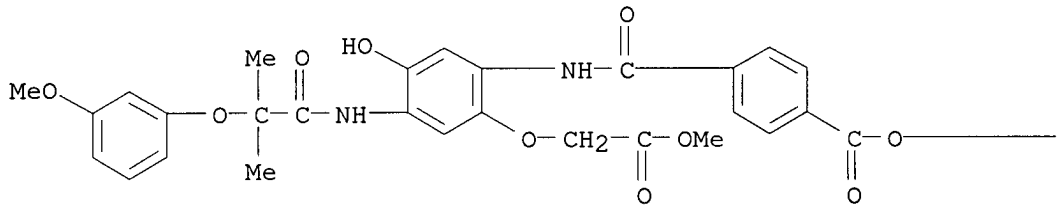

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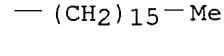
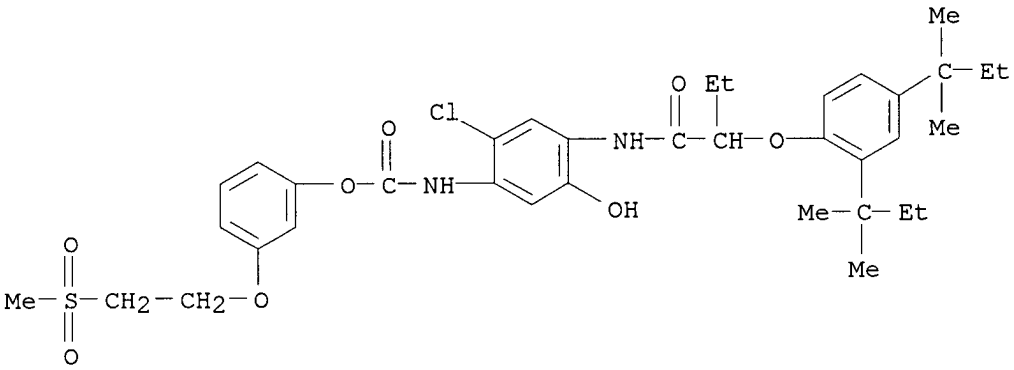


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

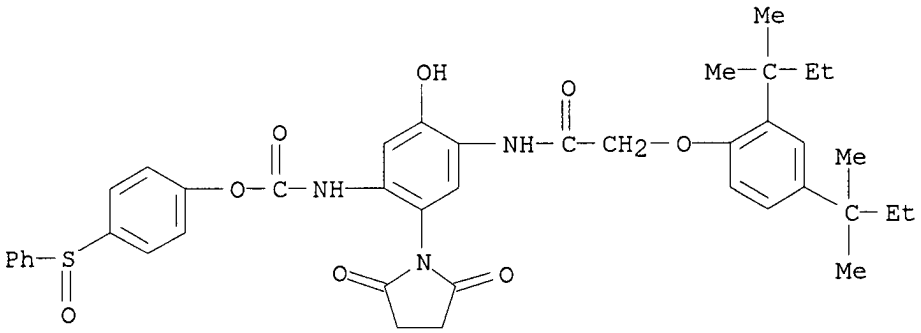
L3 ANSWER 37 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Carbamic acid, [4-[[2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-1-oxobutyl]amino]-2-chloro-5-hydroxyphenyl]-, 3-[2-(methylsulfonyl)ethoxy]phenyl ester (9CI)
MF C36 H47 Cl N2 O8 S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

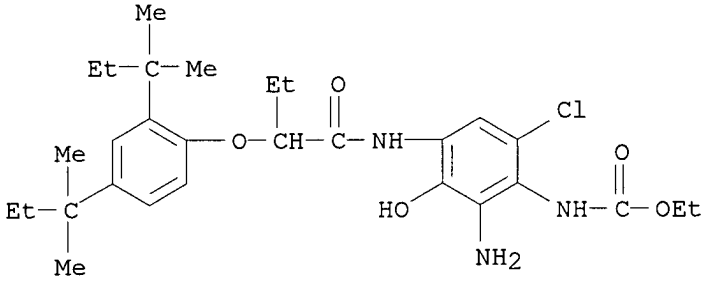
L3 ANSWER 38 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Carbamic acid, [4-[[[2,4-bis(1,1-dimethylpropyl)phenoxy]acetyl]amino]-2-

(2,5-dioxo-1-pyrrolidinyl)-5-hydroxyphenyl]-, 4-(phenylsulfinyl)phenyl ester (9CI)
MF C41 H45 N3 O8 S



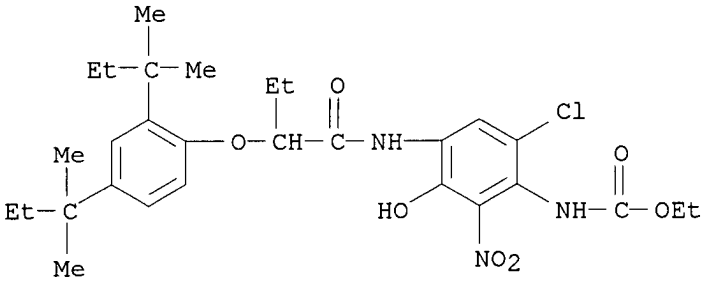
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 39 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Carbamic acid, [2-amino-4-[[2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-1-oxobutyl]amino]-6-chloro-3-hydroxyphenyl]-, ethyl ester (9CI)
MF C29 H42 Cl N3 O5



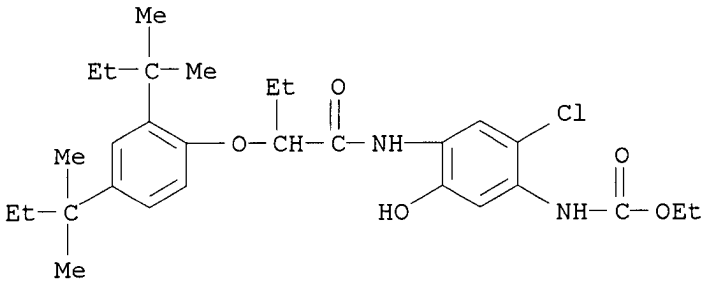
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 40 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Carbamic acid, [4-[[2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-1-oxobutyl]amino]-6-chloro-3-hydroxy-2-nitrophenyl]-, ethyl ester (9CI)
MF C29 H40 Cl N3 O7



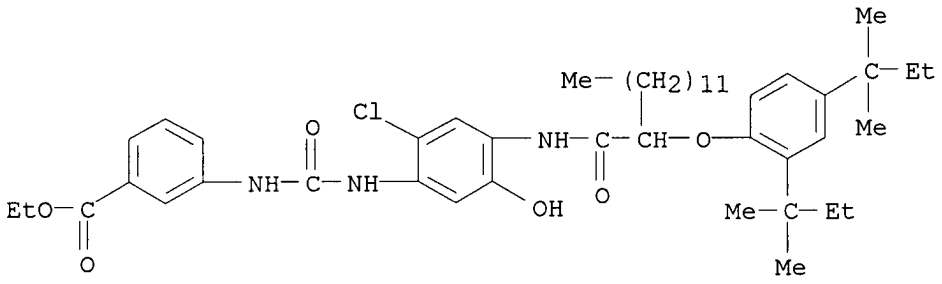
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 41 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Carbamic acid, [4-[[2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-1-oxobutyl]amino]-2-chloro-5-hydroxyphenyl]-, ethyl ester (9CI)
MF C29 H41 Cl N2 O5



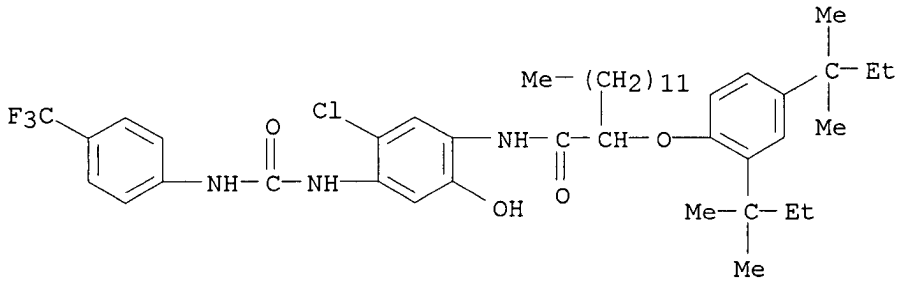
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 42 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Benzoic acid, 3-[[[[4-[[2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-1-oxotetradecyl]amino]-2-chloro-5-hydroxyphenyl]amino]carbonyl]amino]-, ethyl ester (9CI)
MF C46 H66 Cl N3 O6



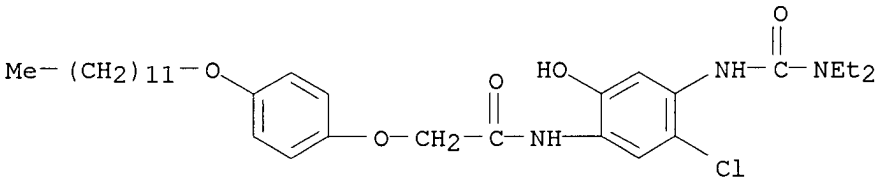
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 43 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Tetradecanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[5-chloro-2-hydroxy-4-[[[[4-(trifluoromethyl)phenyl]amino]carbonyl]amino]phenyl]- (9CI)
MF C44 H61 Cl F3 N3 O4



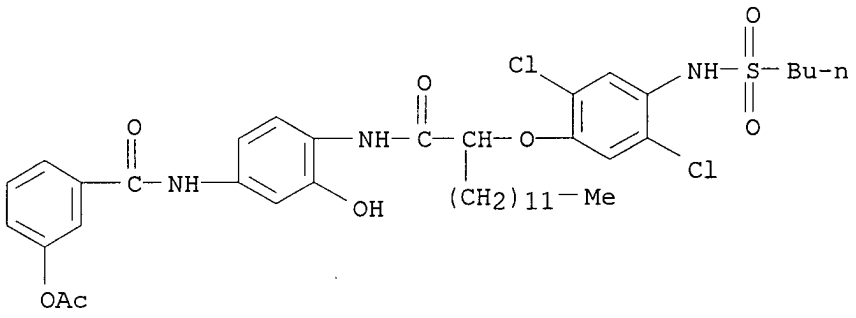
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 44 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Acetamide, N-[5-chloro-4-[[(diethylamino)carbonyl]amino]-2-hydroxyphenyl]-2-[4-(dodecyloxy)phenoxy]- (9CI)
MF C31 H46 Cl N3 O5



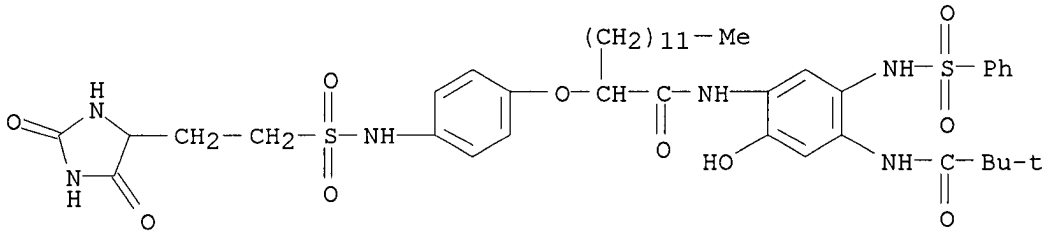
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 45 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Benzamide, 3-(acetyloxy)-N-[4-[[2-[4-[(butylsulfonyl)amino]-2,5-dichlorophenoxy]-1-oxotetradecyl]amino]-3-hydroxyphenyl]- (9CI)
MF C39 H51 Cl2 N3 O8 S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 46 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Tetradecanamide, N-[4-[(2,2-dimethyl-1-oxopropyl)amino]-2-hydroxy-5-[(phenylsulfonyl)amino]phenyl]-2-[4-[[[2-(2,5-dioxo-4-imidazolidinyl)ethyl]sulfonyl]amino]phenoxy]- (9CI)
MF C42 H58 N6 O10 S2



L3 ANSWER 47 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Benzoic acid, 4-[1-[[[4-[(2,2,3,3,4,4,4-heptafluoro-1-oxopropyl)amino]-2-hydroxyphenyl]amino]carbonyl]propoxy]-, 2-(dodecylphenylamino)-2-oxo-1-phenylethyl ester (9CI)
MF C47 H52 F7 N3 O7

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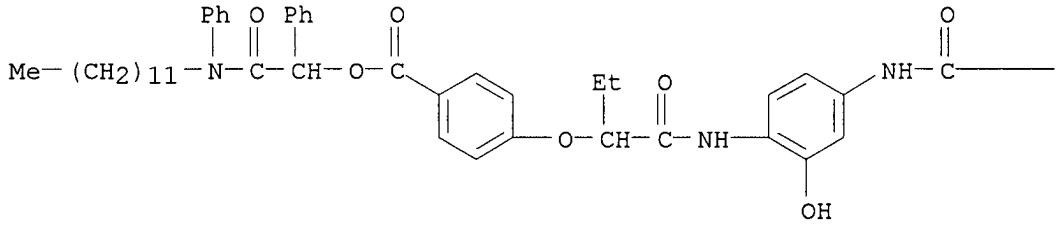

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—CF_2—CF_2—CF_3

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 48 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Benzoic acid, 4-[1-[[[4-[[2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-1-oxohexyl]amino]-2-hydroxyphenyl]amino]carbonyl]propoxy]-, 2-(octadecylphenylamino)-2-oxo-1,1-diphenylethyl ester (9CI)
MF C77 H103 N3 O8

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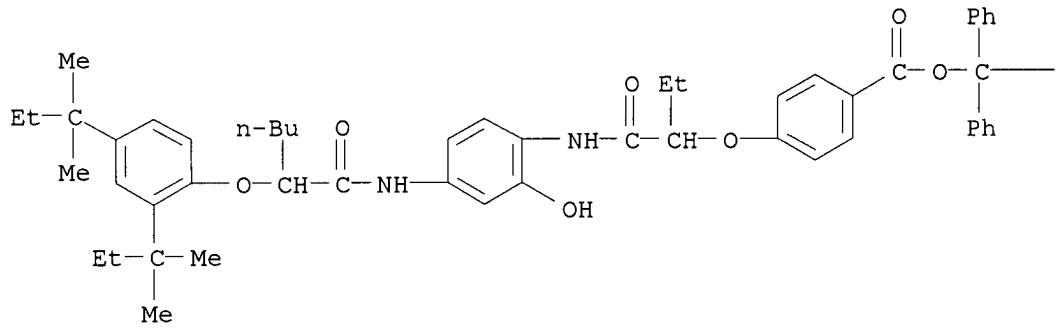

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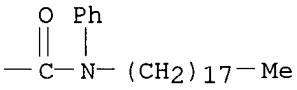

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 49 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Benzoic acid, 4,4'-[(2-hydroxy-1,4-phenylene)bis[imino(2-oxo-2,1-ethanediyl)oxy]]bis-, bis[2-[[4-(diethylamino)phenyl]amino]-2-oxo-1-phenylethyl] ester (9CI)
MF C60 H60 N6 O11

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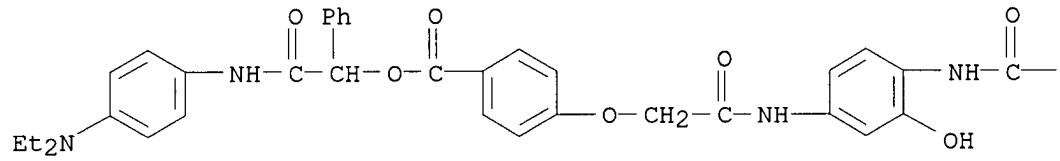

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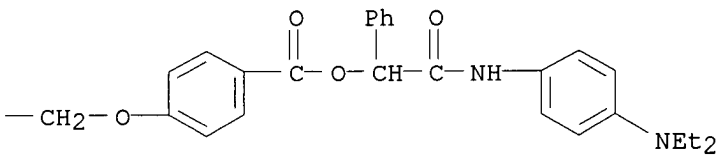

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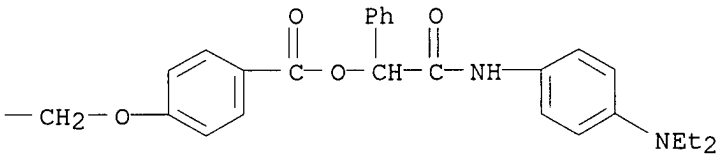

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 50 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Benzoic acid, 4-[2-[[4-[[2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-1-oxobutyl]amino]-2-hydroxyphenyl]amino]-2-oxoethoxy]-, 1,1-dimethyl-2-[[4-(4-morpholinyl)phenyl]amino]-2-oxoethyl ester (9CI)
MF C49 H62 N4 O9

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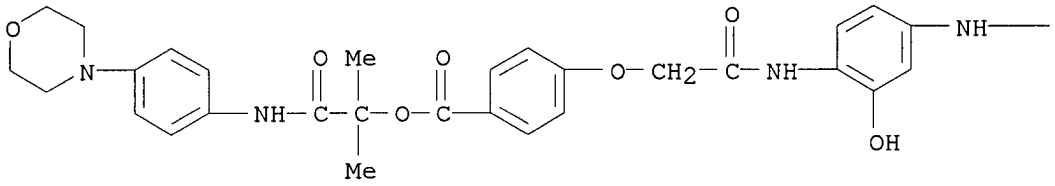

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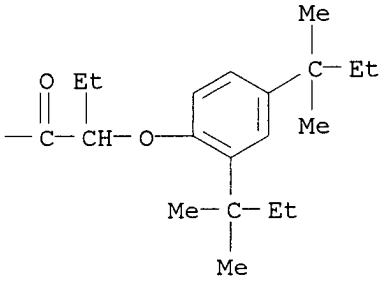
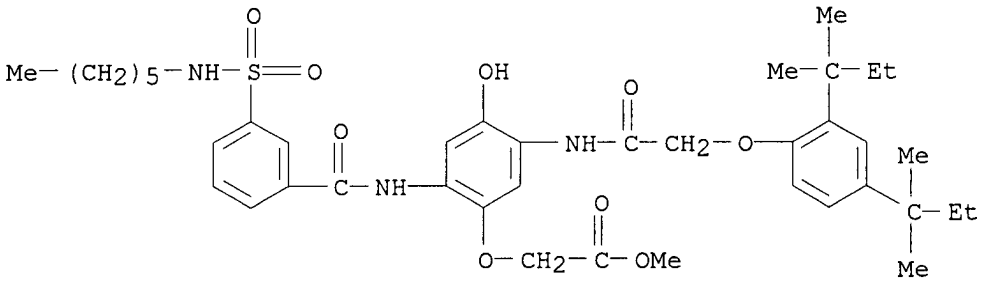


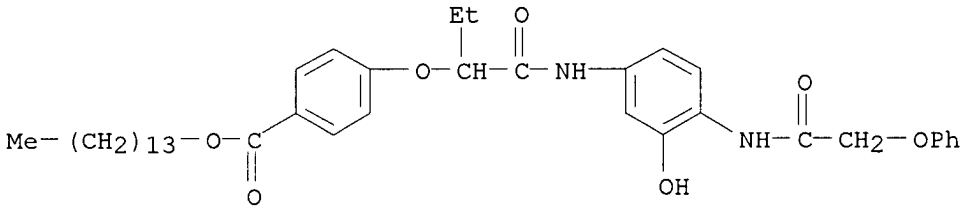
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 51 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Acetic acid, [5-[[[2,4-bis(1,1-dimethylpropyl)phenoxy]acetyl]amino]-2-[[3-[(hexylamino)sulfonyl]benzoyl]amino]-4-hydroxyphenoxy]-, methyl ester (9CI)
MF C40 H55 N3 O9 S



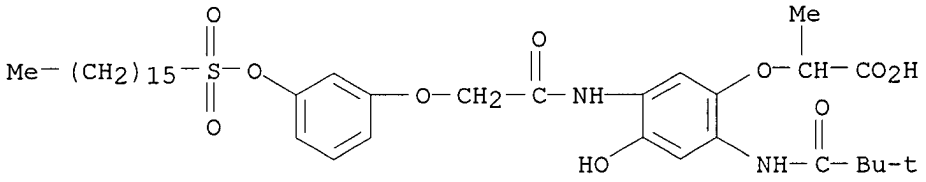
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 52 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Benzoic acid, 4-[1-[[[3-hydroxy-4-[(phenoxyacetyl)amino]phenyl]amino]carbonyl]propoxy]-, tetradecyl ester (9CI)
MF C39 H52 N2 O7



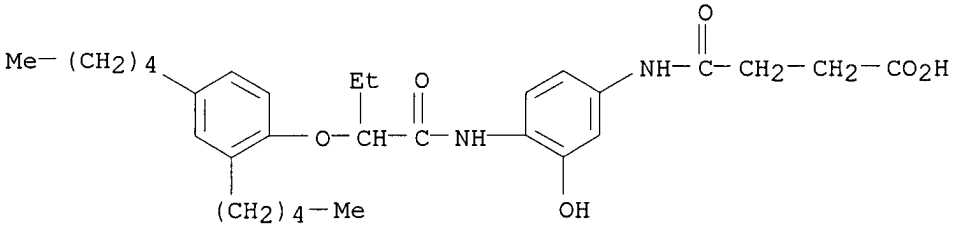
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 53 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Propanoic acid, 2-[2-[(2,2-dimethyl-1-oxopropyl)amino]-5-[[[3-[(hexadecylsulfonyl)oxy]phenoxy]acetyl]amino]-4-hydroxyphenoxy]- (9CI)
MF C38 H58 N2 O10 S



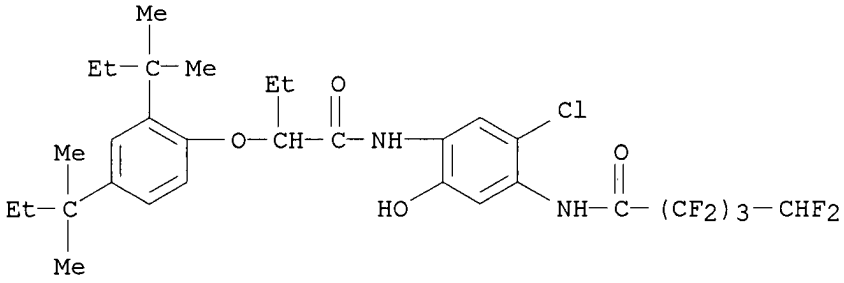
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 54 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Butanoic acid, 4-[[4-[[2-(2,4-dipentylphenoxy)-1-oxobutyl]amino]-3-hydroxyphenyl]amino]-4-oxo- (9CI)
MF C30 H42 N2 O6



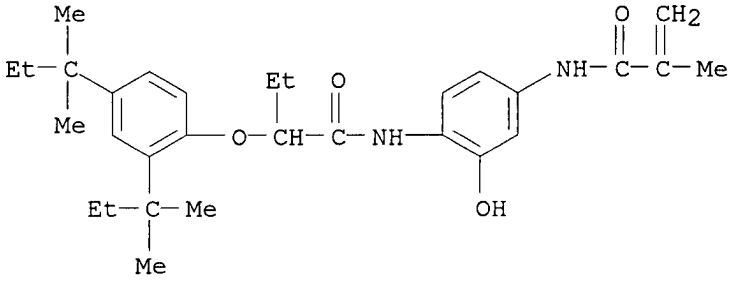
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 55 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Pentanamide, N-[4-[[2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-1-oxobutyl]amino]-2-chloro-5-hydroxyphenyl]-2,2,3,3,4,4,5,5-octafluoro- (9CI)
MF C31 H37 Cl F8 N2 O4



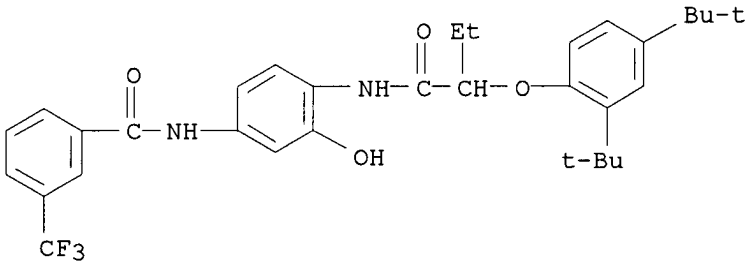
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 56 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN Butanamide, 2-[2,4-bis(1,1-dimethylpropyl)phenoxy]-N-[2-hydroxy-4-[(2-methyl-1-oxo-2-propenyl)amino]phenyl]- (9CI)
MF C30 H42 N2 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 ANSWER 57 OF 57 REGISTRY COPYRIGHT 2002 ACS
IN m-Toluanilide, 4'-[2-(2,4-di-tert-butylphenoxy)butyramido]-
    .alpha.,.alpha.,.alpha.-trifluoro-3'-hydroxy- (7CI, 8CI)
MF C32 H37 F3 N2 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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Uploading 09856396.str

L4 STRUCTURE UPLOADED

=> s l4 full
FULL SEARCH INITIATED 09:03:04 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 2532 TO ITERATE

100.0% PROCESSED 2532 ITERATIONS 1910 ANSWERS
SEARCH TIME: 00.00.01

L5 1910 SEA SSS FUL L4

=>
Uploading 09856396.str

L6 STRUCTURE UPLOADED

=> s l6 full
FULL SEARCH INITIATED 09:05:52 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 456 TO ITERATE

100.0% PROCESSED 456 ITERATIONS 2 ANSWERS
SEARCH TIME: 00.00.01

L7 2 SEA SSS FUL L6

=> d 1-2 bib ab
'BIB' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'
'AB' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual
fields or predefined formats. The predefined substance formats
are: (RN = CAS Registry Number)

REG - RN
SAM - Index Name, MF, and structure - no RN
FIDE - All substance data, except sequence data
IDE - FIDE, but only 50 names
SQIDE - IDE, plus sequence data
SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used
SQD - Protein sequence data, includes RN
SQD3 - Same as SQD, but 3-letter amino acid codes are used
SQN - Protein sequence name information, includes RN

CALC - Table of calculated properties
EPROP - Table of experimental properties
PROP - EPROP and CALC

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract
APPS -- Application and Priority Information
BIB -- CA Accession Number, plus Bibliographic Data
CAN -- CA Accession Number
CBIB -- CA Accession Number, plus Bibliographic Data (compressed)
IND -- Index Data
IPC -- International Patent Classification
PATS -- PI, SO
STD -- BIB, IPC, and NCL

IABS --ABS, indented, with text labels
IBIB -- BIB, indented, with text labels
ISTD -- STD format, indented

OBIB -------- AN, plus Bibliographic Data (original)
OIBIB ------- OBIB, indented with text labels

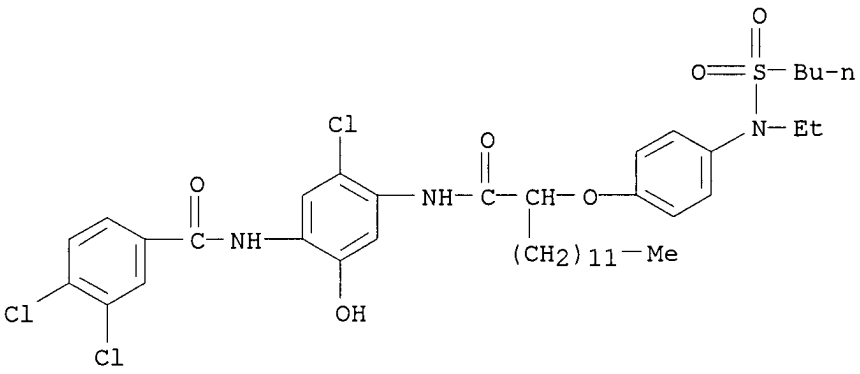
SBIB -------- BIB, no citations
SIBIB ------- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.
The MAX format is the same as ALL.
The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

For additional information, please consult the following help messages:

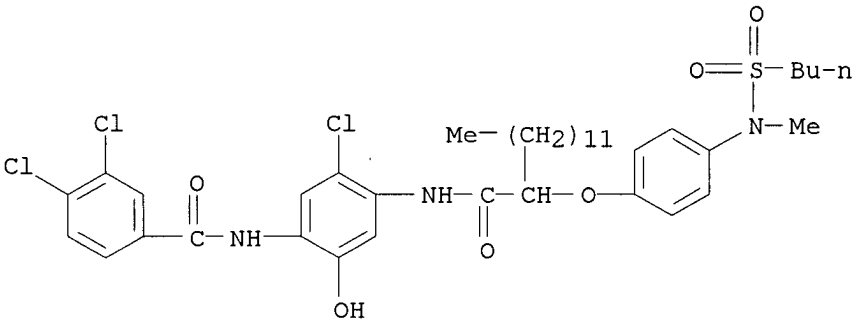
HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
ENTER DISPLAY FORMAT (IDE):sam

L7 ANSWER 1 OF 2 REGISTRY COPYRIGHT 2002 ACS
IN Benzamide, N-[4-[[2-[4-[(butylsulfonyl)ethylamino]phenoxy]-1-oxotetradecyl]amino]-5-chloro-2-hydroxyphenyl]-3,4-dichloro- (9CI)
MF C39 H52 Cl3 N3 O6 S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 ANSWER 2 OF 2 REGISTRY COPYRIGHT 2002 ACS
IN Benzamide, N-[4-[[2-[4-[(butylsulfonyl)methylamino]phenoxy]-1-oxotetradecyl]amino]-5-chloro-2-hydroxyphenyl]-3,4-dichloro- (9CI)
MF C38 H50 Cl3 N3 O6 S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	471.80	472.01

FILE 'CAPLUS' ENTERED AT 09:06:19 ON 29 OCT 2002
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FILE COVERS 1907 - 29 Oct 2002 VOL 137 ISS 18
FILE LAST UPDATED: 27 Oct 2002 (20021027/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

CAS roles have been modified effective December 16, 2001. Please check your SDI profiles to see if they need to be revised. For information on CAS roles, enter HELP ROLES at an arrow prompt or use the CAS Roles thesaurus (/RL field) in this file.

=> d his

(FILE 'HOME' ENTERED AT 08:56:18 ON 29 OCT 2002)

FILE 'REGISTRY' ENTERED AT 08:56:34 ON 29 OCT 2002
L1 STRUCTURE UPLOADED
L2 4 S L1 SSS
L3 57 S L1 FULL
L4 STRUCTURE UPLOADED
L5 1910 S L4 FULL
L6 STRUCTURE UPLOADED
L7 2 S L6 FULL

FILE 'CAPLUS' ENTERED AT 09:06:19 ON 29 OCT 2002

=> s l7
L8 1 L7

=> d bib ab

L8 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2002 ACS
AN 2000:442031 CAPLUS
DN 133:65924
TI Photographic coupler providing dye images with improved light stability
IN Clarke, David; Leyshon, Llewellyn J.; Smith, Katie E.
PA Eastman Kodak Company, USA
SO PCT Int. Appl., 42 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2000038013	A1	20000629	WO 1999-GB4360	19991222
	W: JP, US				
	RW: AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE				
	EP 1145078	A1	20011017	EP 1999-962405	19991222
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, FI				
	JP 2002533760	T2	20021008	JP 2000-590010	19991222
PRAI	GB 1998-28147	A	19981222		
	WO 1999-GB4360	W	19991222		

OS MARPAT 133:65924
AB There is provided a photog. element comprising at least one silver halide emulsion layer having assocd. therewith a novel phenolic cyan dye-forming photog. coupler of formula I [X = H or a group that can be split off by

the reaction of the coupler with an oxidized color developing agent; Y, Z = (substituted) alkyl, (substituted) aryl, a (substituted) 5-10-membered heterocyclic ring which contains one or more heteroatoms selected from N, O, and S, or II where R, R2 = (substituted) alkyl, (substituted) aryl, or a (substituted) 5-10-membered heterocyclic ring which contains one or more heteroatoms selected from N, O, and S; R1, R3 = H, (substituted) alkyl, or (substituted) aryl; n = 1 or 2]. The photog. coupler provides cyan dye images having a good hue and an excellent resistance to heat, humidity, and light.

RE.CNT 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> d his

(FILE 'HOME' ENTERED AT 08:56:18 ON 29 OCT 2002)

FILE 'REGISTRY' ENTERED AT 08:56:34 ON 29 OCT 2002
L1 STRUCTURE UPLOADED
L2 4 S L1 SSS
L3 57 S L1 FULL
L4 STRUCTURE UPLOADED
L5 1910 S L4 FULL
L6 STRUCTURE UPLOADED
L7 2 S L6 FULL

FILE 'CAPLUS' ENTERED AT 09:06:19 ON 29 OCT 2002
L8 1 S L7